

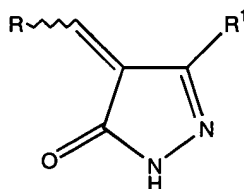
Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

Claims 1-17 (Cancelled)

18. (Currently Amended) A compound represented by the following structural formula:



or physiologically acceptable salts thereof, wherein:

R is selected from the group consisting of substituted or unsubstituted pyrrolyl;

wherein R can be substituted by one or more halogen, lower alkyl group, R³O-, hydroxyl, HOC(O), R³OC(O)-, R³OC(O)R⁶-, R³OR⁶-, trihalomethyl, trihalomethylcarbonyl, nitro, -C(O)NR⁴R⁵, -NR⁴R⁵, R³CO-, -(CH₂)ₙR⁷, -C(O)(CH₂)ₙ-R⁷, -C(O)(CH₂)ₙ-C(O)-R⁷, -O(CH₂)ₙR⁷, -C(O)NR⁴(CH₂)ₙR⁷, -C(O)O(CH₂)ₙR⁷, -OC(O)(CH₂)ₙR⁷, -NR⁴C(O)(CH₂)ₙR⁷, -R⁶NR⁴R⁵, -R⁶N(R⁴)-R⁶-R⁷, -R⁶N(R⁶-R⁷)₂, -R⁶C(O)NR⁴(CH₂)ₙR⁷, -R⁶C(O)O(CH₂)ₙR⁷, -R⁶OC(O)(CH₂)ₙR⁷, -R⁶NR⁴C(O)(CH₂)ₙR⁷, -R⁶CH(C(O)OR⁴)(NR⁵C(O)R⁴) or a substituted aryl or aralkyl group, wherein the substituent is selected from the group consisting of halogen, trihalomethyl, hydroxy, -NR⁴R⁵, nitro, -CONR⁴R⁵, lower alkyl group, R³O-, -C(O)OR⁴ or -OC(O)R³;

wherein R⁶ is a lower alkyl group or an aryl group;

wherein R⁷ is alkoxy, haloalkyl, lower alkyl piperazine, hydroxyl, R³O-, R³C(O)- or -NR⁴R⁵;

wherein suitable substituents for R³, R⁴ and R⁵ can be one or more moieties selected from the group consisting of halogens, lower alkyl, hydroxy, lower alkoxy, carboxy, lower alkyl ester, trihalomethyl, nitro, phenyl, phenyl-lower alkyl, (C₃-C₆)cycloalkyl, (C₃-C₆)cycloalkyl-alkyl, CN, amino, alkylamino, dialkylamino, -C(O)NH₂, -C(O)NH(alkyl) and -C(O)N(alkyl)₂;

R¹ is hydrogen or -A-Z;

A is -(CH₂)ₙ;

Z is a ring system selected from the group consisting of phenyl, pyrazinyl, wherein said ring system can be optionally substituted with one or more moieties selected from the group consisting of halogens, lower alkyl, R^3O- , $HO-$, $HOC(O)-$, $R^3OC(O)-$, trihalomethyl, nitro, an aromatic group, a (C_3-C_6) cycloalkyl group, a heterocyclic group, an aralkyl group, a (C_3-C_6) cycloalkyl-alkyl group, a heterocycl-alkyl group, $-CN$, $-C(O)NR^4R^5$ or $-NR^4R^5$;

R^3 for each occurrence is, independently selected from the group consisting of substituted or unsubstituted: lower alkyl group, lower alkoxy lower alkyl group, aromatic group, (C_3-C_6) cycloalkyl group, heterocyclic group, aralkyl group, a (C_3-C_6) cycloalkyl-alkyl group, and heterocycl-alkyl group;

R^4 and R^5 for each occurrence are each, independently, hydrogen, or are selected from the group consisting of substituted or unsubstituted: lower alkyl group, aromatic group, (C_3-C_6) cycloalkyl group, heterocyclic group, aralkyl group, a (C_3-C_6) cycloalkyl-alkyl group, and heterocycl-alkyl group;

optionally, R^4 and R^5 together with the nitrogen to which they are attached represent morpholino, pyrrolidino, piperidino, imidazol-1-yl, piperazino, thiamorpholino, azepino or perhydro-1,4-diazepin-1-yl groups each optionally substituted by one or more moieties selected from the group consisting of lower alkyl, hydroxy, lower alkoxy lower alkyl, an aromatic group, a (C_3-C_6) cycloalkyl group, a heterocyclic group, an aralkyl group, a (C_3-C_6) cycloalkyl-alkyl group, and a heterocycl-alkyl group; and

n is 0.

Claim 19-22 (Cancelled)

23. (Original) The compound of Claim 18 wherein the compound is a mixture of stereoisomers.

24. (Original) The compound of Claim 23 wherein the stereoisomers are enantiomers.

25. (Original) The compound of Claim 24 wherein the stereoisomers are E and Z isomers.

26. (Original) The compound of claim 18 wherein the compound is a mixture of structural isomers.

27. (Original) The compound of claim 26 wherein the structural isomers are tautomers.

Claims 28 - 40 (Cancelled)

41. (Currently Amended) A compound according to Claim 18 wherein R is substituted with one or more substituents, each independently selected from the group consisting of halogens, lower alkyl groups, R^3O- , hydroxyl, $HOC(O)$, $R^3OC(O)-$, $R^3OC(O)R^6-$, R^3OR^6- , trihalomethyl, trihalomethylcarbonyl, nitro, $-C(O)NR^4R^5$, $-NR^4R^5$, R^3CO- , $(CH_2)_n-R^7$, $-C(O)(CH_2)_nR^7$, $-O(CH_2)_nR^7$, $-C(O)NR^4(CH_2)_nR^7$, $-C(O)O(CH_2)_nR^7$, $-OC(O)(CH_2)_nR^7$, $-NR^4C(O)(CH_2)_nR^7$, $-R^6NR^4R^5$, $-R^6N(R^4)-R^6-R^7$, $-R^6N(R^6-R^7)_2$, $-R^6C(O)NR^4(CH_2)_nR^7$, $-R^6C(O)O(CH_2)_nR^7$, $-R^6OC(O)(CH_2)_nR^7$, $-R^6NR^4C(O)(CH_2)_nR^7$, $-R^6CH(C(O)OR^4)(\overset{NR^5}{C}NR^5C(O)R^4)$, an optionally substituted aryl and an optionally substituted aralkyl group;

wherein the optionally substituted aryl and optionally substituted aralkyl groups are optionally substituted with one or more substituents selected from the group consisting of halogen, trihalomethyl, hydroxyl, $-NR^4R^5$, nitro, $-CONR^4R^5$, lower alkyl group, R^3O- , $-C(O)OR^4$ and $-OC(O)R^3$;

R^6 is a lower alkyl group or an aryl group; and

R^7 is alkoxy, haloalkyl, loweralkyl piperazine, hydroxyl, R^3O- , $R^3C(O)-$ or $-NR^4R^5$.

42. (Cancelled)

43. (Previously Presented) A compound of claim 41, wherein R is pyrrol-2-yl or pyrrol-3-yl.

44. (Cancelled)

45. (Previously Presented) A compound of Claim 43, wherein R is optionally substituted with one or more moieties selected from the group consisting of Br, Cl, F, aminomethyl, N,N-dimethylaminomethyl, carboxy, carboxymethyl, carboxyethyl, carbonylmethyl, carbonylethyl, methoxycarbonyl, ethoxycarbonyl, phenyl, 4-morpholinomethyl, $-C(O)-O-(CH_2)_2-N(Me)_2$, $-C(O)-O-(CH_2)_2-N(Et)_2$, $-C(O)-O-CH_2-N(Me)_2$, $-C(O)-O-(CH_2)_2-N(Me)_2$, $-C(O)-NH-(CH_2)_2-N(Me)_2$, $-CH_2-NH-C(O)-CF_3$, $(CH_2)_n-R^7$ and an optionally substituted moiety selected from the group consisting of methyl, ethyl, propyl, isopropyl, butyl and phenyl, where said optionally substituted

moiety is optionally substituted with one or more of Br, Cl, F, hydroxyl, nitro, amino or lower alkyl.

46. (Previously Presented) A compound of claim 18 wherein R¹ is pyrazinyl or phenyl and R is pyrrolyl substituted by one or more methyl and diethylaminoethyl.

47. (Previously Presented) A compound of claim 46 wherein the compound is 4-[4-(2-diethylamino-ethyl)-3,5-dimethyl-1H-pyrrol-2-ylmethylene]-5-pyrazin-2-yl-2,4-dihydropyrazol-3-one.